=>	d	hi	

(FILE 'HOME' ENTERED AT 23:51:23 ON 27 APR 2006)

FILE 'STNGUIDE' ENTERED AT 23:51:58 ON 27 APR 2006

FILE 'REGISTRY' ENTERED AT 23:59:58 ON 27 APR 2006

L11 S MYCOPHENOLIC ACID/CN

FILE 'CAPLUS, BIOSIS, USPATFULL, DRUGU' ENTERED AT 00:02:15 ON 28 APR 2006

L2 3015 S L1

L3 31 S POLYMER AND BACKBONE AND ANHYDRIDE LINKAGE AND L2 AND HYDROLY

L431 DUP REMOVE L3 (0 DUPLICATES REMOVED)

=>

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L1 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN RN 24280-93-1 REGISTRY
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ED Entered STN: 16 Nov 1984

CN 4-Hexenoic acid, 6-(1,3-dihydro-4-hydroxy-6-methoxy-7-methyl-3-oxo-5-isobenzofuranyl)-4-methyl-, (4E)- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN 4-Hexenoic acid, 6-(1,3-dihydro-4-hydroxy-6-methoxy-7-methyl-3-oxo-5-isobenzofuranyl)-4-methyl-, (E)-

CN 4-Hexenoic acid, 6-(4-hydroxy-6-methoxy-7-methyl-3-oxo-5-phthalanyl)-4-methyl-, (E)- (8CI)

OTHER NAMES:

CN Lilly 68618

CN Melbex

CN MPA

CN Mycophenolic acid

CN NSC 129185

FS STEREOSEARCH

MF C17 H20 O6

CI COM

LC STN Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IMSPATENTS, IMSRESEARCH, IPA, MEDLINE, MRCK*, NAPRALERT, NIOSHTIC, PATDPASPC, PHAR, PROMT, PS, RTECS*, SCISEARCH, SYNTHLINE, TOXCENTER, USAN, USPAT2, USPATFULL

(*File contains numerically searchable property data)

Other Sources: EINECS**, WHO

(**Enter CHEMLIST File for up-to-date regulatory information)

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1192 REFERENCES IN FILE CA (1907 TO DATE)

71 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

1195 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> file caplus biosis uspatf drugu COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 8.42 9.41

FULL ESTIMATED COST

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L1 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN
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RN 24280-93-1 REGISTRY

ED Entered STN: 16 Nov 1984

CN 4-Hexenoic acid, 6-(1,3-dihydro-4-hydroxy-6-methoxy-7-methyl-3-oxo-5-isobenzofuranyl)-4-methyl-, (4E)- (9CI) (CA INDEX NAME)

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OTHER NAMES:

CN Lilly 68618

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FS STEREOSEARCH

MF C17 H20 O6

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Other Sources: EINECS**, WHO

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Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1192 REFERENCES IN FILE CA (1907 TO DATE)

71 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

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